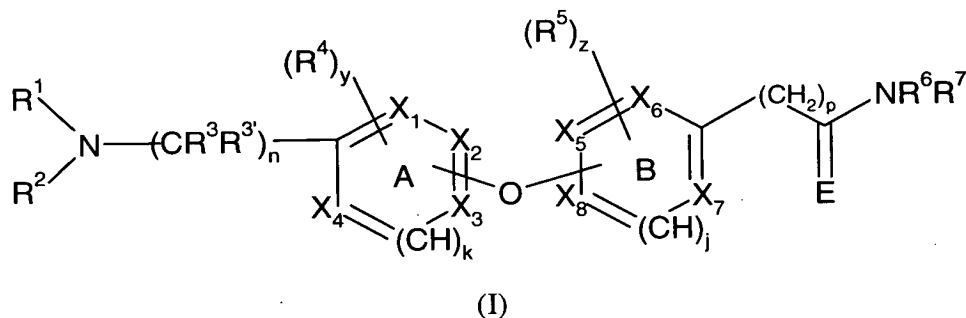


Amendments to the Claims

Listing of Claims:

1. (Original) A compound of the formula (I)



wherein

when K is 0, one of X₁, X₂, X₃, X₄, is a S or O atom and the others are independently selected from C, CH, or N; and wherein when j is 0, one of X₅, X₆, X₇, and X₈ is S, or O, and the others are independently selected from C, CH, or N; provided that both k and j are not simultaneously equal to zero or 1; and provided that each of rings A or B has no more than 2 nitrogen atoms;

n is 0, 1, 2, or 3;

k is 0 or 1; j is 0 or 1;

p is 0, 1 or 2;

E is O or NH;

R¹ and R² are independently selected from hydrogen, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, phenyl, C₁-C₁₀ alkylaryl, C(O)C₁-C₈ alkyl, CO(O)C₁-C₈alkyl, SO₂C₁-C₈ alkyl, SO₂C₁-C₁₀ alkylaryl, or SO₂C₁-C₈ alkylheterocyclic, C₄-C₁₀ alkylcycloalkane, C₁-C₈ alkoxyalkyl, (CH₂)_nC(O)OR⁸, (CH₂)_nC(O)R⁸, (CH₂)_mC(O)NR⁸R⁸, and (CH₂)_mNSO₂R⁸; wherein each of the alkyl, alkenyl, and aryl groups are optionally substituted with one to five groups independently selected from C₁-C₈ alkyl, C₂-C₈ alkenyl, phenyl, C₁-C₈ alkylaryl, C(O)C₁-C₈ alkyl, CO(O)C₁-C₈ alkyl, C₁-C₈ alkoxy, SO₂C₁-C₈ alkyl, SO₂C₁-C₈ alkylaryl, SO₂C₁-C₈ alkylheterocyclic, C₄-C₁₀ alkylcycloalkane, (CH₂)_nC(O)OR⁸, (CH₂)_nC(O)R⁸; and wherein R¹ and R² may optionally combine with each other to form a 4, 5, 6, or 7-member nitrogen-containing heterocycle which nitrogen-containing heterocycle may have substituents

selected from the group consisting of oxo, amino, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, phenyl, C₁-C₈ alkylaryl, C(O)C₁-C₈ alkyl, CO(O)C₁-C₈ alkyl, halo, C₁-C₈ haloalkyl;

R³ and R^{3'} are each independently selected from Hydrogen, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, C₁-C₈ alkoxy, C₁-C₈ thioalkyl, phenyl, aryl, C₁-C₈ alkylaryl;

R⁴ and R⁵ are each independently selected from Hydrogen, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, C₁-C₈ alkoxy, halo, C₁-C₈ haloalkyl, phenyl, aryl, C₁-C₈ alkylaryl, (CH₂)_mNSO₂C₁-C₈ alkyl, (CH₂)_mNSO₂phenyl, (CH₂)_mNSO₂aryl, -C(O)C₁-C₈ alkyl, or -C(O)OC₁-C₈ alkyl; wherein each R⁴ or R⁵ is attached to its respective ring only at carbon atoms, and wherein y is 0, 1, 2, or 3; and wherein z is 0, 1, 2, or 3;

R⁶ and R⁷ are each independently selected from hydrogen, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, C(O)C₁-C₈ alkyl, hydroxy, C₁-C₈ alkoxy, SO₂C₁-C₈ alkyl, SO₂C₁-C₈ alkylaryl, or SO₂C₁-C₈ alkylheterocyclic, aryl, C₁-C₈ alkylaryl, C₃-C₇ cycloalkane, C₁-C₁₀ alkylcycloalkane, (CH₂)_nC(O)OR⁸, (CH₂)_nC(O)R⁸, (CH₂)_mC(O)NR⁸R⁸, and (CH₂)_mNSO₂R⁸; wherein each of the alkyl, alkenyl, and aryl groups are optionally substituted with one to five groups independently selected from C₁-C₈ alkyl, C₂-C₈ alkenyl, phenyl, and C₁-C₈ alkylaryl; and wherein R⁶ and R⁷ may independently combine together, and with the nitrogen atom to which they are attached or with 0, 1, or 2 atoms adjacent to the nitrogen atom to which they are attached to form a 4, 5, 6, or 7-membered nitrogen containing heterocycle which nitrogen containing heterocycle may further have substituents selected from the group consisting of oxo, amino, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₂-C₈ alkynyl, phenyl, C₁-C₈ alkylaryl, C(O)C₁-C₈ alkyl, CO(O)C₁-C₈ alkyl, hydroxy, C₁-C₈ alkoxy, halo, and haloalkyl;

R⁸ is hydrogen, C₁-C₈ alkyl, C₂-C₈ alkenyl, C₅-C₈ alkylaryl, -C(O)C₁-C₈ alkyl, or -C(O)OC₁-C₈ alkyl; m is 1, 2, or 3; or a pharmaceutically acceptable salt, solvate, enantiomer, racemate, diastereomers or mixtures thereof.

2. (Original) The compound according to claim 1 wherein the A-ring is selected from the group consisting of phenyl, pyridyl, thiophene, thiazole, furanyl, imidazole, and pyrazole.

3. (Cancelled)

4. (Original) A compound according to Claim 1 wherein the B-ring is selected from the group consisting of phenyl, thiophene, thiazole, furan, and pyridine.

5. (Cancelled)

6. (Cancelled)

7. (Cancelled)

8. (Cancelled)

9. (Cancelled)

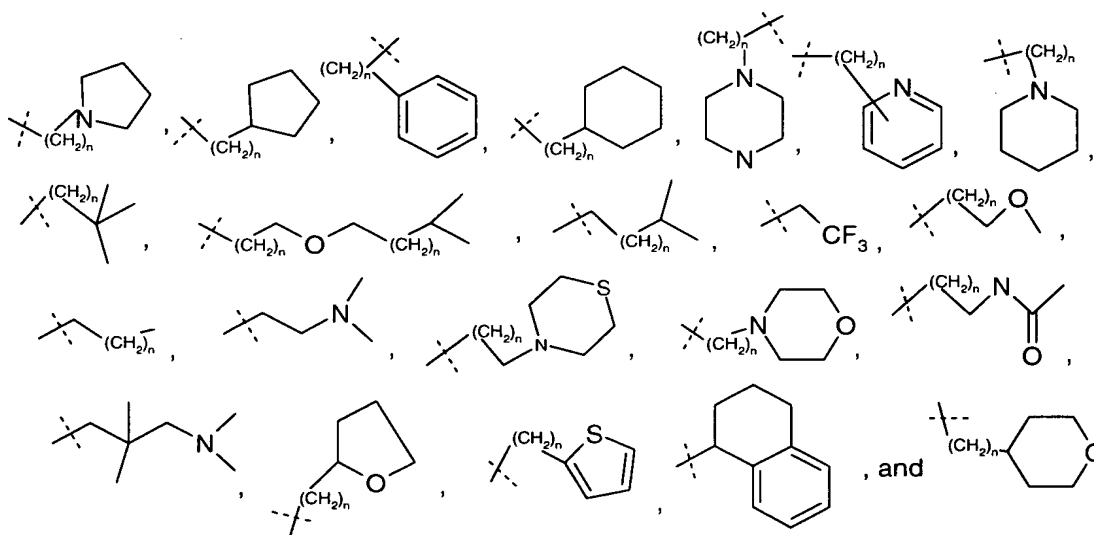
10. (Cancelled)

11. (Cancelled)

12. (Original) A compound according to Claim 1 wherein y is 0, or 1, and R⁴ is independently selected from the group consisting of hydrogen, fluoro, chloro, methyl, methoxy, ethoxy, ethyl, isopropyl, trifluoromethyl, phenyl, and benzyl.

13. (Original) A compound according to Claim 1 wherein z is 0, or 1, and R⁵ is independently selected from the group consisting of hydrogen, fluoro, chloro, methyl, methoxy, ethoxy, ethyl, isopropyl, trifluoromethyl, phenyl, and benzyl.

14. (Original) A compound according to Claim 1 wherein R¹ and R² are each independently selected from the group consisting of hydrogen, methyl, ethyl, propyl, isopropyl, phenyl,



15. (Original) The compound according to Claim 1 wherein R⁶ and R⁷ are each independently selected from the group consisting of hydrogen, methyl, ethyl, propyl, isopropyl, and phenyl.

16. (Cancelled)

17. (Cancelled)

18. (Cancelled)

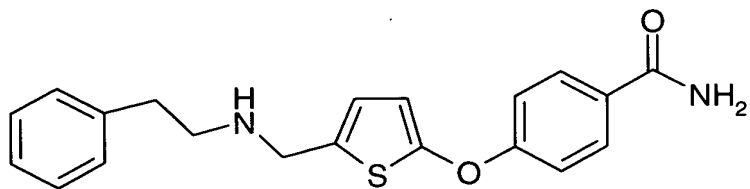
19. (Cancelled)

20. (Original) A compound selected from the group consisting of:
5-{4-[(3-Methyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{4-[(3,3-Dimethyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{4-[(2-Cyclopentyl-ethylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{4-[(3-Ethyl-pentylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{4-[(Cyclohexylmethyl-amino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-(4-{[2-(4-Fluoro-phenyl)-ethylamino]-methyl}-phenoxy)-thiophene-2-carboxamide,
5-{2-Chloro-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{2-Chloro-4-[(3,3-dimethyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-(2-Chloro-4-{[2-(4-fluoro-phenyl)-ethylamino]-methyl}-phenoxy)-thiophene-2-carboxamide,
5-{2-Fluoro-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-{4-[(3,3-Dimethyl-butylamino)-methyl]-2-fluoro-phenoxy}-thiophene-2-carboxamide,
5-(2-Fluoro-4-{[2-(4-fluoro-phenyl)-ethylamino]-methyl}-phenoxy)-thiophene-2-carboxamide,
5-{2-Methoxy-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiophene-2-carboxamide,
5-(4-{[2-(4-Fluoro-phenyl)-ethylamino]-methyl}-2-methoxy-phenoxy)-thiophene-2-carboxamide,
4-{5-[(3-Methyl-butylamino)-methyl]-thiazol-2-yloxy}-benzamide,
3-Methoxy-4-{5-[(3-methyl-butylamino)-methyl]-thiazol-2-yloxy}-benzamide,

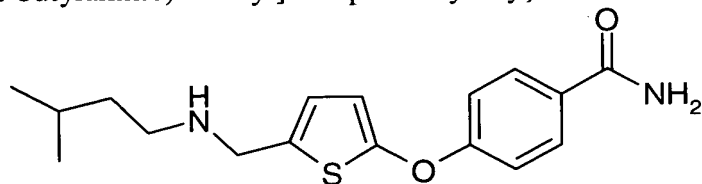
4-{5-[(3,3-Dimethyl-butylamino)-methyl]-thiazol-2-yloxy}-benzamide,
 4-(5-{[2-(4-Fluoro-phenyl)-ethylamino]-methyl}-thiazol-2-yloxy)-benzamide,
 4-(5-{[2-(4-Fluoro-phenyl)-ethylamino]-methyl}-thiazol-2-yloxy)-3-methoxy-
 benzamide,
 4-{5-[(Cyclohexylmethyl-amino)-methyl]-thiazol-2-yloxy}-benzamide,
 2-(4-Pentylaminomethyl-phenoxy)-thiazole-5-carboxamide,
 2-{4-[(3-Methyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 2-{4-[(3,3-Dimethyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 2-(4-{[2-(4-Fluoro-phenyl)-ethylamino]-methyl}-phenoxy)-thiazole-5-carboxamide,
 2-{2-Chloro-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 2-{2-Fluoro-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 2-{2-Methyl-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 2-{2-Methoxy-4-[(3-methyl-butylamino)-methyl]-phenoxy}-thiazole-5-carboxamide,
 4-[5-(2,6-Dimethyl-morpholin-4-ylmethyl)-thiazol-2-yloxy]-benzamide,
 4-{5-[(3-Methoxy-propylamino)-methyl]-thiazol-2-yloxy}-benzamide,
 4-{4-Chloro-5-[(3-methyl-butylamino)-methyl]-thiazol-2-yloxy}-benzamide,
 4-(5-Butylaminomethyl-4-chloro-thiazol-2-yloxy)-benzamide,
 4-{4-Chloro-5-[(3,3-dimethyl-butylamino)-methyl]-thiazol-2-yloxy}-benzamide,
 4-[5-(Phenethylamino-methyl)-thiophen-2-yloxy]-benzamide,
 4-{5-[(3-Methyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide,
 4-(5-{[2-(3-Fluoro-phenyl)-ethylamino]-methyl}-thiophen-2-yloxy)-benzamide,
 4-{5-[(2-Cyclopentyl)-ethylamino]-methyl}-thiophen-2-yloxy}-benzamide,
 4-{5-[(2-Thiophen-2-yl)-ethylamino]-methyl}-thiophen-2-yloxy}-benzamide,
 4-{5-[(3,3-Dimethyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide,
 3-Methoxy-4-[5-(phenethylamino-methyl)-thiophen-2-yloxy]-benzamide
 hydrochloride,
 3-Methoxy-4-{5-[(3-methyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide
 hydrochloride,
 4-[5-(2-Phenethylamino-ethyl)-thiophen-2-yloxy]-benzamide hydrochloride,
 or a pharmaceutically acceptable salt, solvate, enantiomer, diastereomer or
 diastereomeric mixture thereof.

21. (Original) A compound selected from the group consisting of:

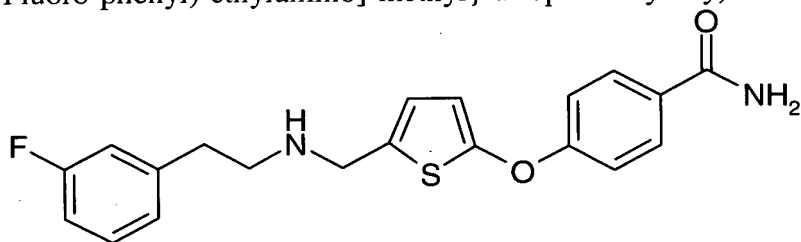
4-[5-(Phenethylamino-methyl)-thiophen-2-yloxy]-benzamide



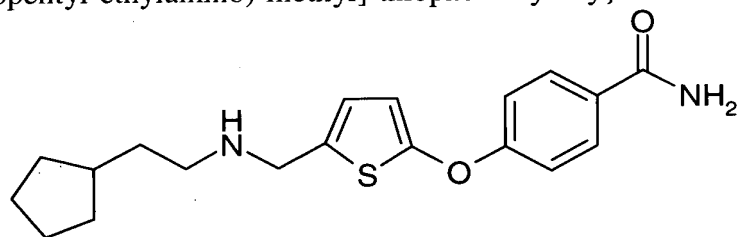
4-{5-[(3-Methyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide



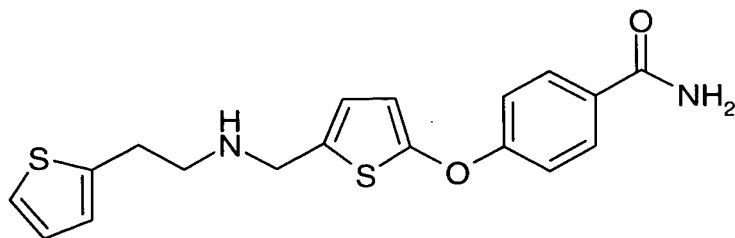
4-(5-{[2-(3-Fluoro-phenyl)-ethylamino]-methyl}-thiophen-2-yloxy)-benzamide



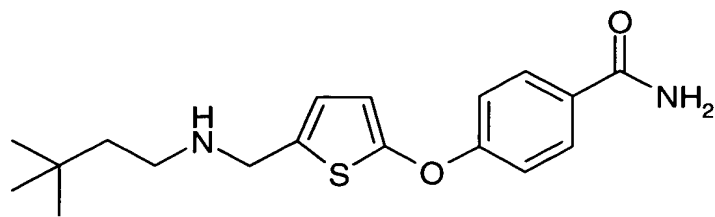
4-{5-[(2-Cyclopentyl-ethylamino)-methyl]-thiophen-2-yloxy}-benzamide



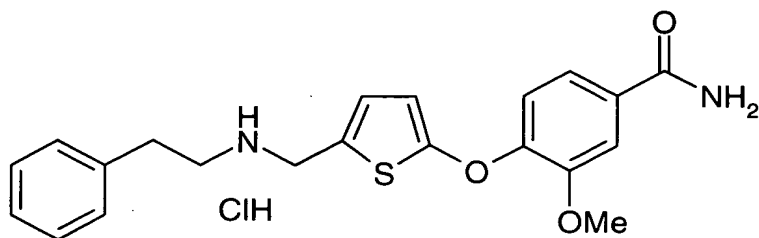
4-{5-[(2-Thiophen-2-yl-ethylamino)-methyl]-thiophen-2-yloxy}-benzamide



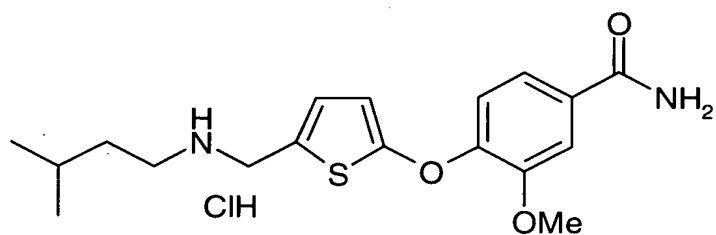
4-{5-[(3,3-Dimethyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide



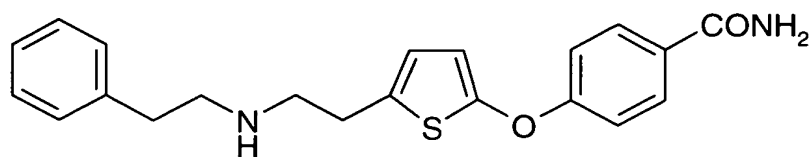
3-Methoxy-4-[5-(phenethylamino-methyl)-thiophen-2-yloxy]-benzamide hydrochloride



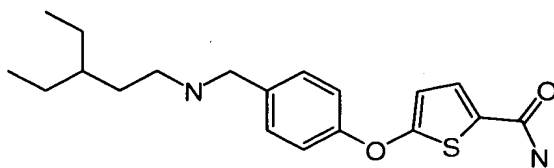
3-Methoxy-4-{5-[(3-methyl-butylamino)-methyl]-thiophen-2-yloxy}-benzamide hydrochloride



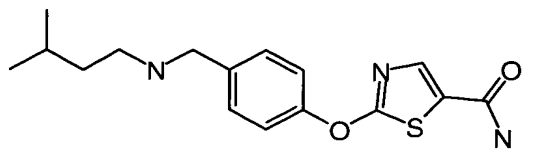
4-[5-(2-Phenethylamino-ethyl)-thiophen-2-yloxy]-benzamide hydrochloride



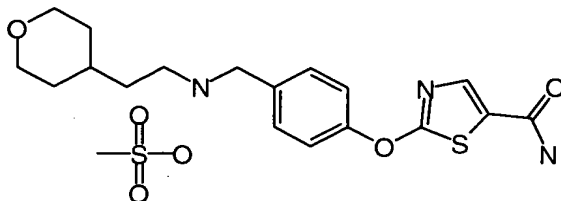
5-{4-[(3-Ethylpentylamino)methyl]phenoxy}thiophene-2-carboxamide



2-{4-[(3-Methylbutylamino)methyl]phenoxy}thiazole-5-carboxamide



2-(4-([2-(Tetrahydropyran-4-yl)ethylamino] methyl)phenoxy)thiazole-5-carboxamide methanesulfonate



and a pharmaceutically acceptable salt, solvate, enantiomer, diastereomer or diastereomeric mixture thereof.

22. (Original) A pharmaceutical composition comprising a therapeutically effective amount of a compound of formula I in association with a carrier, diluent and/or excipient.

23. (Cancelled)

24. (Cancelled)

25. (Cancelled)

26. (Currently Amended) A method ~~according to Claim 25 wherein the~~ treating or preventing obesity and Related Diseases is selected from the group consisting of diabetes, diabetic complications, diabetic retinopathy, atherosclerosis, hyperlipidemia, hypertriglycemia, hyperglycemia, and hyperlipoproteinemia, irritable bowel syndrome, depression, smoking and alcohol addiction, sexual dysfunction, substance abuse, drug overdose, addictive behavior disorders, compulsive behaviors, and stroke comprising administering a therapeutically effective amount of a compound of formula I.

27. (Original) A method of suppressing appetite comprising administering a therapeutically effective amount of a compound of formula I to a patient in need thereof.

28. (Cancelled)